W O R L D W I D E

wwPDB

PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

May 29, 2020 – 05:09 am BST

PDB ID	:	4F2N
Title	:	Crystal structure of iron superoxide dismutase from Leishmania major
Authors	:	Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on	:	2012-05-08
Resolution	:	1.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

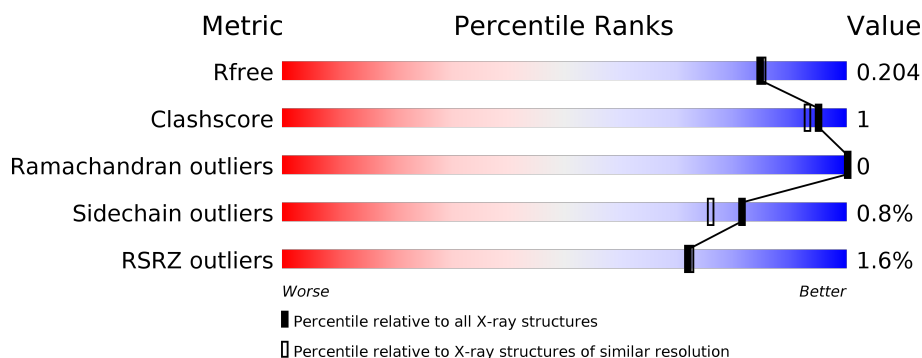|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.

Metric	Percentile Ranks	Value
Rfree		0.204
Clashscore		1
Ramachandran outliers		0
Sidechain outliers		0.8%
RSRZ outliers		1.6%

Worse   *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	230	2% 87% • 9%
1	B	230	87% • 9%
1	C	230	% 87% • 9%
1	D	230	86% 5% 9%
1	E	230	% 86% 5% 9%
1	F	230	% 90% • 9%

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Mol	Chain	Length	Quality of chain
1	G	230	2% ... 87% ... • ... 9%
1	H	230	% ... 86% ... 5% ... 9%
1	I	230	% ... 87% ... • ... 9%
1	J	230	% ... 86% ... 5% ... 9%
1	K	230	3% ... 83% ... 7% • ... 9%
1	L	230	3% ... 90% ... • ... 9%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 21527 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Superoxide dismutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total 1683	C 1082	N 281	O 310	S 10	0	2	0
1	B	210	Total 1682	C 1082	N 283	O 307	S 10	0	1	0
1	C	209	Total 1667	C 1072	N 280	O 305	S 10	0	1	0
1	D	210	Total 1691	C 1090	N 283	O 308	S 10	0	1	0
1	E	209	Total 1659	C 1067	N 276	O 306	S 10	0	0	0
1	F	210	Total 1698	C 1093	N 285	O 310	S 10	0	3	0
1	G	209	Total 1679	C 1079	N 281	O 309	S 10	0	2	0
1	H	210	Total 1682	C 1082	N 283	O 307	S 10	0	1	0
1	I	209	Total 1665	C 1072	N 280	O 303	S 10	0	0	0
1	J	210	Total 1675	C 1080	N 280	O 305	S 10	0	0	0
1	K	209	Total 1655	C 1065	N 276	O 304	S 10	0	0	0
1	L	209	Total 1657	C 1067	N 278	O 302	S 10	0	0	0

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total 1	Fe 1	0	0
2	J	1	Total 1	Fe 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	Fe 1	0	0
2	K	1	Total 1	Fe 1	0	0
2	E	1	Total 1	Fe 1	0	0
2	H	1	Total 1	Fe 1	0	0
2	B	1	Total 1	Fe 1	0	0
2	I	1	Total 1	Fe 1	0	0
2	C	1	Total 1	Fe 1	0	0
2	A	1	Total 1	Fe 1	0	0
2	L	1	Total 1	Fe 1	0	0
2	F	1	Total 1	Fe 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	167	Total 167	O 167	0	0
3	B	159	Total 159	O 159	0	0
3	C	120	Total 120	O 120	0	0
3	D	122	Total 122	O 122	0	0
3	E	124	Total 124	O 124	0	0
3	F	103	Total 103	O 103	0	0
3	G	129	Total 129	O 129	0	0
3	H	112	Total 112	O 112	0	0
3	I	92	Total 92	O 92	0	0

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| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	J	144	Total 144	O 144	0	0
3	K	77	Total 77	O 77	0	0
3	L	73	Total 73	O 73	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.
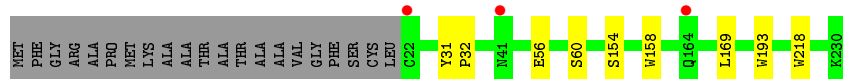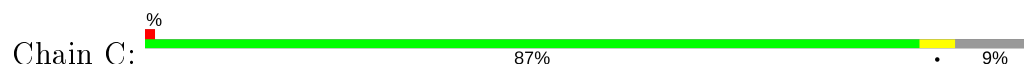
- Molecule 1: Superoxide dismutase
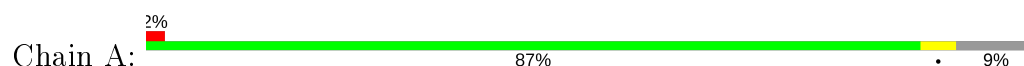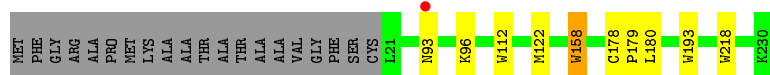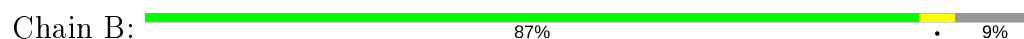
- Molecule 1: Superoxide dismutase

- Molecule 1: Superoxide dismutase

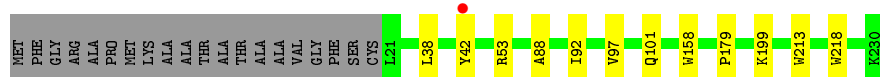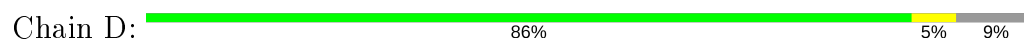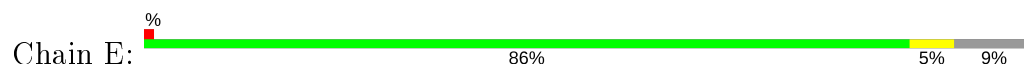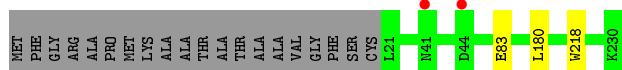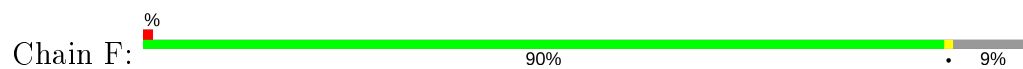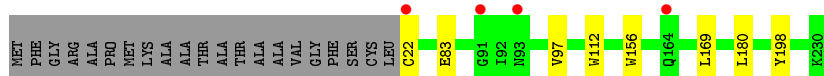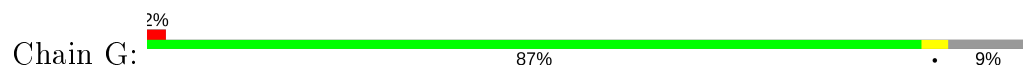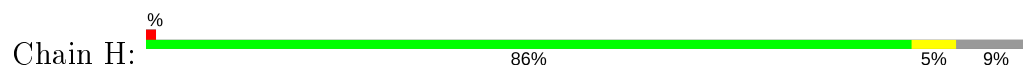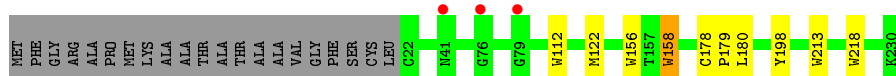- Molecule 1: Superoxide dismutase

- Molecule 1: Superoxide dismutase

- Molecule 1: Superoxide dismutase

Chain F:

- Molecule 1: Superoxide dismutase

Chain G:

- Molecule 1: Superoxide dismutase

Chain H:

- Molecule 1: Superoxide dismutase

Chain I:

- Molecule 1: Superoxide dismutase

Chain J:

- Molecule 1: Superoxide dismutase

Chain K:

- Molecule 1: Superoxide dismutase

Chain L:
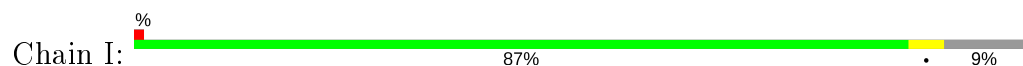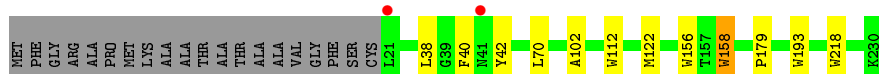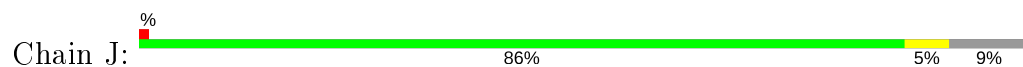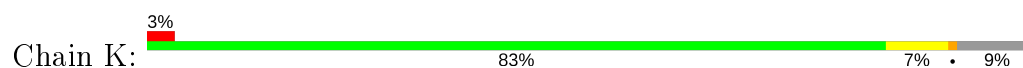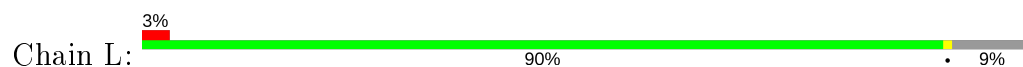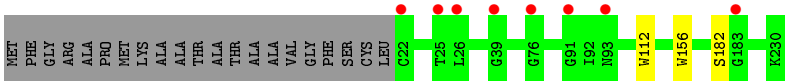
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	47.03Å 102.74Å 168.24Å 91.66° 96.36° 95.01°	Depositor
Resolution (Å)	19.77 – 1.85 19.77 – 1.85	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.77-1.85) 96.3 (19.77-1.85)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	2.66 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.173 , 0.203 0.175 , 0.204	Depositor DCC
R_{free} test set	12890 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	25.3	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 44.1	EDS
L-test for twinning[2]	$<|L|> = 0.48$, $<L^2> = 0.31$	Xtriage
Estimated twinning fraction	0.008 for -h,-k,h+l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	21527	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $<|L|>$, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: FE2

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$|Z|$>5	RMSZ	#$|Z|$>5
1	A	0.71	1/1742 (0.1%)	0.65	0/2370
1	B	0.75	3/1737 (0.2%)	0.65	0/2361
1	C	0.65	2/1722 (0.1%)	0.60	0/2342
1	D	0.69	2/1748 (0.1%)	0.61	0/2376
1	E	0.71	2/1712 (0.1%)	0.68	0/2332
1	F	0.66	1/1760 (0.1%)	0.63	0/2392
1	G	0.72	2/1739 (0.1%)	0.63	0/2368
1	H	0.70	5/1738 (0.3%)	0.62	0/2363
1	I	0.68	4/1717 (0.2%)	0.62	0/2333
1	J	0.70	4/1728 (0.2%)	0.63	0/2350
1	K	0.63	3/1708 (0.2%)	0.59	0/2327
1	L	0.63	2/1709 (0.1%)	0.59	0/2324
All	All	0.69	31/20760 (0.1%)	0.62	0/28238

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	156	TRP	CD2-CE2	6.63	1.49	1.41
1	J	193	TRP	CD2-CE2	6.17	1.48	1.41
1	E	156	TRP	CD2-CE2	6.09	1.48	1.41
1	I	112	TRP	CD2-CE2	5.81	1.48	1.41
1	J	158	TRP	CD2-CE2	5.78	1.48	1.41

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1683	0	1587	5	0
1	B	1682	0	1595	6	0
1	C	1667	0	1569	3	0
1	D	1691	0	1601	5	0
1	E	1659	0	1542	3	0
1	F	1698	0	1615	1	0
1	G	1679	0	1567	2	0
1	H	1682	0	1590	5	0
1	I	1665	0	1577	3	0
1	J	1675	0	1582	5	0
1	K	1655	0	1538	7	0
1	L	1657	0	1560	0	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	167	0	0	0	0
3	B	159	0	0	1	0
3	C	120	0	0	0	0
3	D	122	0	0	0	0
3	E	124	0	0	0	0
3	F	103	0	0	0	1
3	G	129	0	0	1	0
3	H	112	0	0	0	0
3	I	92	0	0	0	0
3	J	144	0	0	0	0
3	K	77	0	0	1	0
3	L	73	0	0	0	1
All	All	21527	0	18923	45	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

The worst 5 of 45 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:38:LEU:HD12	1:K:42:TYR:HB3	1.68	0.76
1:A:38:LEU:HD12	1:A:42:TYR:HB3	1.76	0.66
1:D:38:LEU:HD12	1:D:42:TYR:HB3	1.87	0.56
1:J:38:LEU:HD12	1:J:42:TYR:HB3	1.87	0.55
1:B:178:CYS:SG	1:B:180:LEU:HG	2.46	0.54

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:489:HOH:O	3:L:470:HOH:O[1_545]	2.11	0.09

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	209/230 (91%)	204 (98%)	5 (2%)	0	100	100
1	B	209/230 (91%)	203 (97%)	6 (3%)	0	100	100
1	C	208/230 (90%)	204 (98%)	4 (2%)	0	100	100
1	D	209/230 (91%)	205 (98%)	4 (2%)	0	100	100
1	E	207/230 (90%)	202 (98%)	5 (2%)	0	100	100
1	F	211/230 (92%)	206 (98%)	5 (2%)	0	100	100
1	G	209/230 (91%)	205 (98%)	4 (2%)	0	100	100
1	H	209/230 (91%)	202 (97%)	7 (3%)	0	100	100
1	I	207/230 (90%)	204 (99%)	3 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	J	208/230 (90%)	203 (98%)	5 (2%)	0	100	100
1	K	207/230 (90%)	204 (99%)	3 (1%)	0	100	100
1	L	207/230 (90%)	203 (98%)	4 (2%)	0	100	100
All	All	2500/2760 (91%)	2445 (98%)	55 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	179/195 (92%)	177 (99%)	2 (1%)	73	65
1	B	178/195 (91%)	178 (100%)	0	100	100
1	C	175/195 (90%)	174 (99%)	1 (1%)	86	83
1	D	179/195 (92%)	179 (100%)	0	100	100
1	E	173/195 (89%)	169 (98%)	4 (2%)	50	34
1	F	181/195 (93%)	181 (100%)	0	100	100
1	G	177/195 (91%)	174 (98%)	3 (2%)	60	47
1	H	178/195 (91%)	177 (99%)	1 (1%)	86	83
1	I	175/195 (90%)	174 (99%)	1 (1%)	86	83
1	J	176/195 (90%)	176 (100%)	0	100	100
1	K	172/195 (88%)	169 (98%)	3 (2%)	60	47
1	L	173/195 (89%)	172 (99%)	1 (1%)	86	83
All	All	2116/2340 (90%)	2100 (99%)	16 (1%)	81	76

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	22	CYS
1	G	169	LEU

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Mol	Chain	Res	Type
1	K	169	LEU
1	E	182	SER
1	K	198	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	164	GLN
1	J	24	HIS

5.3.3   RNA (i)

There are no RNA molecules in this entry.

5.4   Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5   Carbohydrates (i)

There are no carbohydrates in this entry.

5.6   Ligand geometry (i)

Of 12 ligands modelled in this entry, 12 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7   Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	209/230 (90%)	-0.38	4 (1%)	66	66	14, 23, 36, 57	0
1	B	210/230 (91%)	-0.32	1 (0%)	91	91	15, 24, 37, 50	0
1	C	209/230 (90%)	-0.27	3 (1%)	75	76	18, 29, 41, 54	0
1	D	210/230 (91%)	-0.23	1 (0%)	91	91	18, 27, 43, 61	0
1	E	209/230 (90%)	-0.19	3 (1%)	75	76	14, 26, 44, 60	1 (0%)
1	F	210/230 (91%)	-0.21	2 (0%)	82	82	17, 30, 43, 55	0
1	G	209/230 (90%)	-0.28	4 (1%)	66	66	16, 26, 43, 62	0
1	H	210/230 (91%)	-0.25	3 (1%)	75	76	17, 29, 44, 54	0
1	I	209/230 (90%)	-0.18	3 (1%)	75	76	17, 29, 46, 55	0
1	J	210/230 (91%)	-0.29	2 (0%)	82	82	17, 26, 37, 56	0
1	K	209/230 (90%)	-0.09	6 (2%)	51	50	20, 34, 50, 76	0
1	L	209/230 (90%)	0.03	8 (3%)	40	38	20, 35, 54, 99	0
All	All	2513/2760 (91%)	-0.22	40 (1%)	72	72	14, 28, 46, 99	1 (0%)

The worst 5 of 40 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	22	CYS	8.0
1	G	22	CYS	7.2
1	K	22	CYS	6.9
1	E	22	CYS	5.5
1	A	22	CYS	4.8

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($Å^2$)	Q<0.9
2	FE2	D	300	1/1	0.99	0.06	21,21,21,21	0
2	FE2	K	300	1/1	0.99	0.05	21,21,21,21	0
2	FE2	L	300	1/1	0.99	0.04	22,22,22,22	0
2	FE2	I	300	1/1	0.99	0.05	20,20,20,20	0
2	FE2	J	300	1/1	1.00	0.05	19,19,19,19	0
2	FE2	F	300	1/1	1.00	0.07	19,19,19,19	0
2	FE2	H	300	1/1	1.00	0.06	19,19,19,19	0
2	FE2	A	300	1/1	1.00	0.05	17,17,17,17	0
2	FE2	G	300	1/1	1.00	0.05	16,16,16,16	0
2	FE2	C	300	1/1	1.00	0.05	21,21,21,21	0
2	FE2	E	300	1/1	1.00	0.06	16,16,16,16	0
2	FE2	B	300	1/1	1.00	0.07	16,16,16,16	0

6.5 Other polymers (i)

There are no such residues in this entry.